MEMORANDUM

TO: Mr. Addison Rice

Anderson, Mulholland and Associates

DATE: July 2, 2015

FROM: R. Infante

FILE: 1502113AR1

RE:

Data Validation

Air samples

SDG: 1502113AR1

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SUMMARY

Full validation was performed on the data for several gas samples analyzed for selected volatile organic compounds by method Compendium Method TO-15: Determination Of Volatile Organic Compounds (VOCs) In Air Collected In Specially-Prepared Canisters And Analyzed By Gas Chromatography/Mass Spectrometry (GC/MS), January, 1999. The samples were collected at the Bristol Myer Squib-Building 6 VI facility, Humacao, PR site on February 04 and 06, 2015 and submitted to Eurofins Air Toxics, Inc. of Folson, California that analyzed and reported the results under delivery groups (SDG) 1502113AR1.

The sample results were assessed according to USEPA data validation guidance documents in the following order of precedence: Compendium Method TO-15. Determination Of Volatile Organic Compounds (VOCs) In Air Collected In Specially-Prepared Canisters And Analyzed By Gas Chromatography/Mass Spectrometry (GC/MS), January, 1999; Validating Air Samples. Volatile Organic Analysis of Ambient Air in Canisters by Method TO-15, (SOP # HW-31. Revision #4. October, 2006. The QC criteria and data validation actions listed on the data review worksheets are from the primary guidance document, unless otherwise noted. In general the data is valid as reported and may be used for decision making purposes.

The data results are acceptable for use. The following results were qualified as estimated (J): Acetone, Ethanol, and Hexachlorobutadiene in all samples due to the % RSD for the calibration factor or % D in continuing calibration outside method performance criteria and Hexane, Tetrahydrofuran, Cyclohexane, Propylbenzene, Heptane, and 2-Hexanone in samples 1502113AR1-02A/1502113AR1-03A due to the % RSD outside laboratory/method control limit for laboratory duplicates. 2-Propanol concentration was over the calibration range in sample 1502113AR1-04A after dilution. Result qualified as an estimated value (J) and qualified (E) by the laboratory.

SAMPLES The samples included in the review are listed below

Client Sample ID	Lab. Sample ID	Collected Date	Matrix	Analysis
B6-5SS	1502113AR1-01A	02/04/2015	Air	VOCs
B6-3SS	1502113AR1-02A	02/04/2015	Air	VOCs
B6-3SSD	1502113AR1-03A	02/04/2015	Air	VOCs
B6-4SS	1502113AR1-04A	02/06/2015	Air	VOCs

REVIEW ELEMENTS

Sample data were reviewed for the following parameters, where applicable to the method

- Agreement of analysis conducted with chain of custody (COC) form
- o Holding time and sample preservation
- Gas chromatography/mass spectrometry (GC/MS) tunes
- o Initial and continuing calibrations
- o Method blanks/trip blanks/field blank
- o Canister cleaning certification criteria
- Surrogate spike recovery
- o Internal standard performance and retention times
- o Field duplicate results
- o Laboratory control sample/laboratory control sample duplicate (LCS/LCSD) results
- o Quantitation limits and sample results

DISCUSSION

Agreement of Analysis Conducted with COC Request

Sample reports corresponded to the analytical request designated on the chain-of-custody form.

Holding Times and Sample Preservation

Sample preservation was acceptable.

Samples analyzed within method recommended holding time.

GC/MS Tunes

The frequency and abundance of bromofluorobenzene (BFB) tunes were within the QC acceptance criteria. All samples were analyzed within the tuning criteria associated with the method.

Initial and Continuing Calibrations

VOCs (Method TO-15)

The percent relative standard deviations (%RSDs) and response factors (RFs) of all target analytes were within the QC acceptance criteria in the initial calibration. Correlation coefficients (r²) of target analytes were within the QC acceptance criteria. Ongoing accuracy of the instrument was determined by the analysis of a continuing calibration standard. The % RSD for the response factor for acetone and the % D for the ethanol and hexachlorobutadiene continuing calibration response factor outside the method performance criteria:

DATE	LAB FILE	CRITERIA OUT	COMPOUND	SAMPLES
	ID#	RFs, %RSD, <u>%D</u> , r		AFFECTED
02/10/15	v021002.d	- 39 %	Acetone	All samples.
		39 %	Ethanol	
		- 32 %	Hexachlorobutadiene	

Results for acetone, ethanol, and hexachlorobutadiene qualified as estimated (J) in all samples.

Method Blank/Trip Blank/Field Blank

Target analytes were not detected in laboratory method blanks except for the followings:

DATE LAB ID LEVEL/ COMPOUND CONCENTRATION/ ANALYZED MATRIX UNIT

02/10/15 1502113AR1-05A Air/low Acetone 0.32 ppbv

No action taken, 5x acetone concentration in blank < the concentration found in samples.

Summa canister met cleaning certification criteria.

No trip/field blank analyzed with this data package.

Surrogate Spike Recovery

The surrogate recoveries were within the laboratory QC acceptance limits in all samples analyzed.

Internal Standard Performance

VOCs and Methanol (TO-15)

Samples were spiked with the method specified internal standard. Internal standard are performance and retention times met the QC acceptance criteria in all sample analyses and calibration standards.

Laboratory/Field Duplicate Results

Field/laboratory duplicates were analyzed as part of this data set. Target analytes meet the RPD performance criteria of + 25 % for analytes $5 \times SQL$ except for the followings:

Field duplicates:

COMPOUND	SQL	SAMPLE CONC.	DUPLICATE CONC.	RPD	ACTION
Hexane	0.03452	1.2	0.51	81 %	Qualify results (J) in sample and duplicate
Tetrahydrofuran	-	ND	0.26	NR	Qualify results (J) in sample and duplicate
Cyclohexane	0.03469	0.95	0.38	86 %	Qualify results (J) in sample and duplicate
Propylbenzene	0.02205	0.16	0.057	95 %	Qualify results (J) in sample and duplicate
Heptane	0.02709	0.77	0.28	96 %	Qualify results (J) in sample and duplicate
2-Hexanone	-	ND	0.35	93 %	Qualify results (J) in sample and duplicate

LCS/LCSD Results

VOCs

LCS/LCSD (blank spike) associated with this data package were analyzed by the laboratory. Recoveries and RPD within laboratory control limits.

Quantitation Limits and Sample Results

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Dilutions were performed on TO-15 samples (see worksheet). 2-Propanol concentration was over the calibration range in sample 1502113AR1-04A after dilution. Result qualified as an estimated value (J) and qualified (E) by the laboratory.

Calculations were spot checked.

Certification

The following samples 1502113AR1-01A; 1502113AR1-02A; 1502113AR1-03A; and 1502113AR1-04A were analyzed following standard procedures accepted by regulatory agencies. The quality control requirements met the methods criteria except in the occasions described in this document. The results are valid. Some of the results were qualified.

Rafael Infante

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Client Sample ID: B6-5SS Lab ID#: 1502113AR1-01A

MODIFIED EPA METHOD TO-15 GC/MS FULL SCAN

File Name: Dil. Factor:	v021017r1 24.2		of Collection: 2/4 of Analysis: 2/10	
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Freon 12	2.4	Not Detected	12	Not Detected
Freon 114	2.4	Not Detected	17	Not Detected
Chloromethane	12	2.3 J	25	4.7 J
Vinyl Chloride	2.4	Not Detected	6.2	Not Detected
1,3-Butadiene	2.4	Not Detected	5.4	Not Detected
Bromomethane	12	Not Detected	47	Not Detected
Chloroethane	12	Not Detected	32	Not Detected
Freon 11	2.4	Not Detected	14	Not Detected
Ethanol	12	650 J0 🦪	23	1200 J0
Freon 113	2.4	Not Detected	18	Not Detected
1,1-Dichloroethene	2.4	Not Detected	9.6	Not Detected
Acetone	12	150	29	360
2-Propanol	12	390	30	960
Carbon Disulfide	12	6.8 J	38	21 J
3-Chloropropene	12	Not Detected	38	Not Detected
Methylene Chloride	4.8	6.1	17	21
Methyl tert-butyl ether	2.4	210	8.7	760
rans-1,2-Dichloroethene	2.4	Not Detected	9.6	Not Detected
Hexane	2.4	1.5 J	8.5	5.2 J
1,1-Dichloroethane	2.4	Not Detected	9.8	Not Detected
2-Butanone (Methyl Ethyl Ketone)	12	3.4 J	36	10 J
is-1,2-Dichloroethene	2.4	Not Detected	9.6	Not Detected
Tetrahydrofuran	12	3.1 J	36	9.0 J
Chloroform	2.4	6.1	12	30
1,1,1-Trichloroethane	2.4	Not Detected	13	Not Detected
Cyclohexane	2.4	Not Detected	8.3	Not Detected
Carbon Tetrachloride	2.4	Not Detected	15	Not Detected
2,2,4-Trimethylpentane	12	Not Detected	56	Not Detected
Benzene	2.4	1.9 J	7.7	6.1 J
1,2-Dichloroethane	2.4	Not Detected	9.8	Not Detected
Heptane	2.4	93	9.9	380
richloroethene	2.4	Not Detected	9.9 13	Not Detected
1,2-Dichloropropane	2.4	Not Detected	11	Not Detected
1,4-Dioxane	2.4	Not Detected	8.7	Not Detected
Bromodichloromethane	2.4	Not Detected	16	Not Detected
sis-1,3-Dichloropropene	2.4	No.	11	Not Detected
4-Methyl-2-pentanone	2.4	36	9.9	150
Foluene	2.4	880 Infante	9.9	3300
rans-1,3-Dichloropropene	2.4	Not Distended	8 11	Not Detected
1,1,2-Trichloroethane	2.4	Not Detect 4888	S 11	Not Detected
Tetrachloroethene	2.4	Not Detected	<u> </u>	
		NOT Detected	16 50	Not Detected
2-Hexanone	12	New Petected	50	Not Detect



Client Sample ID: B6-5SS Lab ID#: 1502113AR1-01A

MODIFIED EPA METHOD TO-15 GC/MS FULL SCAN

File Name: v021017r1 Date of Collection: 2/4/15 3:49:00 PM
Dil. Factor: 24.2 Date of Analysis: 2/10/15 07:20 PM

Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Dibromochloromethane	2.4	Not Detected	21	Not Detected
1,2-Dibromoethane (EDB)	2.4	Not Detected	18	Not Detected
Chlorobenzene	2.4	Not Detected	11	Not Detected
Ethyl Benzene	2.4	14	10	59
m,p-Xylene	2.4	39	10	170
o-Xylene	2.4	11	10	49
Styrene	2.4	Not Detected	10	Not Detected
Bromoform	2.4	Not Detected	25	Not Detected
Cumene	2.4	0.99 J	12	4.8 J
1,1,2,2-Tetrachloroethane	2.4	Not Detected	17	Not Detected
Propylbenzene	2.4	1.2 J	12	6.0 J
4-Ethyltoluene	2.4	5.2	12	26
1,3,5-Trimethylbenzene	2.4	1.8 J	12	8.6 J
1,2,4-Trimethylbenzene	2.4	4.5	12	22
1,3-Dichlorobenzene	2.4	Not Detected	14	Not Detected
1,4-Dichlorobenzene	2.4	Not Detected	14	Not Detected
alpha-Chlorotoluene	12	Not Detected	63	Not Detected
1,2-Dichlorobenzene	2.4	Not Detected	14	Not Detected
1,2,4-Trichlorobenzene	12	Not Detected	90	Not Detected
Hexachlorobutadiene	12	Not Detected UJ	J 130	Not Detected UJ
Naphthalene	12	Not Detected	63	Not Detected

J = Estimated value.

UJ = Analyte associated with low bias in the CCV and/or LCS. Container Type: 1 Liter Summa Canister (100% Certified)

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	96	70-130
Toluene-d8	99	70-130
4-Bromofluorobenzene	SOCIALO DE	70-130

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J0 = Estimated value due to bias in the CCV.



Client Sample ID: B6-3SS Lab ID#: 1502113AR1-02A

MODIFIED EPA METHOD TO-15 GC/MS FULL SCAN

Compound	File Name: Dil. Factor:	v021016r1 2.29		of Collection: 2/4 of Analysis: 2/10	
Freon 114	Compound	•	Amount	Rpt. Limit	Amount (ug/m3)
Freon 114 0.23 Not Detected 1.6 Not D Chloromethane 1.1 0.72 J 2.4 1. Vinyl Chloride 0.23 Not Detected 0.51 Not D 1,3-Butadiene 0.23 Not Detected 0.51 Not D Bromomethane 1.1 Not Detected 4.4 Not D Chloroethane 1.1 Not Detected 4.4 Not D Freon 11 0.23 0.18 J 1.3 0.1 Ethanol 1.1 16.0 ♂ 2.2 3 Freon 113 0.23 0.059 J 1.8 0.6 7-Freon 113 0.23 Not Detected 0.91 Not D Acetone 1.1 74 ¬ 7 2.7 1 2-Propanol 1.1 58 2.8 1 2-Propanol 1.1 0.21 ¬ 3.6 0.91 3-Chloropropene 1.1 Not Detected 0.91 Not D 4-Propanol 1.1 0.21 ¬	Freon 12	0,23	0.44	1.1	2.2
Chloromethane	Freon 114				Not Detected
Vinyl Chloride 0.23 Not Detected 0.58 Not D J.3-Butadiene 0.23 Not Detected 0.51 Not D Bromomethane 1.1 Not Detected 4.4 Not D Chloroethane 1.1 Not Detected 3.0 Not D Freon 11 0.23 0.18 J 1.3 0.9 Ethanol 1.1 16 J0 J 2.2 30 Freon 113 0.23 0.059 J 1.8 0.4 Acetone 1.1 74 J 2.7 1 Acetone 1.1 74 J 2.7 1 Carbon Disulfide 1.1 0.21 J 3.6 0.6 Carbon Disulfide 1.1 Not Detected 3.6 Not D Carbon Disulfide 1.1 Not Detected 3.6 Not D Carbon Disulfide 1.1 Not Detected 3.6 Not D Methyl tert-butyl ether 0.23 2.1 0.82 7 trans-1,2-Dichloroethane 0.23	Chloromethane	1.1	0.72 J		1.5 J
December 1.1	Vinyl Chloride	0.23	Not Detected		Not Detected
Chloroethane 1.1 Not Detected 3.0 Not D Freon 11 0.23 0.18 J 1.3 0.1 1.1 16 J0 7 2.2 3 0.18 J 1.3 0.1 1.1 16 J0 7 2.2 3 0.059 J 1.8 0.2 1.1 1.1 16 J0 7 2.2 3 0.059 J 1.8 0.2 1.1 1.1 16 J0 7 2.2 3 0.059 J 1.8 0.0 1.1 1.1 16 J0 7 2.2 3 0.059 J 1.8 0.0 1.1 1.1 16 J0 7 2.2 3 0.059 J 1.8 0.0 1.1 1.1 16 J0 7 2.2 7 1 1 0.0 1.1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	1,3-Butadiene	0.23	Not Detected	0.51	Not Detected
Ethanol	Bromomethane	1.1	Not Detected	4.4	Not Detected
Eftenol 11 0.23 0.18 J 1.3 0.8 Ethanol 1.1 16 J0 J 2.2 30 7.1-Dichloroethene 0.23 Not Detected 0.91 Not Detected Acetone 1.1 74 J 2.7 1 2-Propanol 1.1 58 2.8 1 2-Propanol 1.1 58 2.8 1 3-Chloropropene 1.1 Not Detected 3.6 Not Detected 3-Chloropropene 1.1 Not Detected 3.6 Not Detected Methyl tert-butyl ether 0.23 2.1 0.82 7 Trans-1,2-Dichloroethene 0.23 Not Detected 0.91 Not Detected Hexane 0.23 Not Detected 0.91 Not Detected 1,1-Dichloroethane 0.23 Not Detected 0.93 Not Detected 1,1-Dichloroethene 0.23 Not Detected 0.93 Not Detected 1,1-Trichloroethene 0.23 Not Detected 0.91 Not Detected	Chloroethane	1.1	Not Detected	3.0	Not Detected
Freon 113 0.23 0.059 J 1.8 0.4	Freon 11	0.23	0.18 J		0.99 J
1,1-Dichloroethene	Ethanol	1.1	16 J0 3	2.2	30 J0
Acetone 1.1 74 7 2.7 1 2-Propanol 1.1 58 2.8 1 Carbon Disulfide 1.1 0.21 J 3.6 0.6 Carbon Disulfide 1.1 0.21 J 3.6 Not D Methylene Chloride 1.1 Not Detected 3.6 Not D Methylene Chloride 0.46 1.0 1.6 3 Methyl tert-butyl ether 0.23 2.1 0.82 7 Trans-1,2-Dichloroethene 0.23 Not Detected 0.91 Not D 1,1-Dichloroethane 0.23 Not Detected 0.91 Not D 2-Butanone (Methyl Ethyl Ketone) 1.1 4.9 3.4 1 Cis-1,2-Dichloroethene 0.23 Not Detected 0.91 Not D Cis-1,2-Dichloroethene 0.23 Not Detected 0.91 Not D Cis-1,2-Dichloroethane 0.23 Not Detected 0.91 Not D Cis-1,2-Dichloroethane 0.23 Not Detected 0.91 Not D Cis-1,2-Trinichloroethane 0.23 Not Detected 1.2 Not D Cyclohexane 0.23 Not Detected 1.2 Not D Cyclohexane 0.23 0.95 7 0.79 3 Carbon Tetrachloride 0.23 0.95 7 0.79 3 Carbon Tetrachloride 0.23 0.092 J 1.4 0.5 Carbon Tetrachloride 0.093 Not Detected 0.93 Not Detected 0.93 Not Detected 0.93 Not Detected 0.93 Not Dete	Freon 113	0.23	0.059 J	1.8	0.45 J
Acetone 1.1 74 7 2.7 11 2-Propanol 1.1 58 2.8 1 Carbon Disulfide 1.1 0.21 J 3.6 0.6 Acchon Disulfide 1.1 0.21 J 3.6 0.6 Methylene Chloride 1.1 Not Detected 3.6 Not D Methylene Chloride 0.46 1.0 1.6 3 Methyl tert-butyl ether 0.23 2.1 0.82 7 Trans-1,2-Dichloroethene 0.23 Not Detected 0.91 Not D Hexane 0.23 Not Detected 0.91 Not D 2-Butanone (Methyl Ethyl Ketone) 1.1 4.9 3.4 7 Lis-1,2-Dichloroethene 0.23 Not Detected 0.91 Not D Cis-1,2-Dichloroethene 0.23 Not Detected 0.91 Not D Cis-1,2-Dichloroethene 0.23 Not Detected 0.91 Not D Cis-1,2-Dichloroethene 0.23 Not Detected 0.91 Not D Cis-1,2-Trinchloroethane 0.23 Not Detected 0.91 Not D Cyclohexane 0.23 0.067 J 1.1 0.3 1,1,1-Trichloroethane 0.23 Not Detected 1.2 Not D Cyclohexane 0.23 0.095 7 0.79 3 Carbon Tetrachloride 0.23 0.092 J 1.4 0.5 Carbon Tetrachloride 0.23 0.092 J 1.4 0.5 Carbon Tetrachloride 0.23 0.092 J 1.4 0.5 1,2-Dichloroethane 0.23 Not Detected 0.93 Not D Heptane 0.23 Not Detected 0.94 0.65 Romodichloromethane 0.23 Not Detected 1.0 Not D Heptane 0.23 No	1,1-Dichloroethene	0.23	Not Detected	0.91	Not Detected
1.1 58 2.8 1 1.2 3.6 0	•		74 7		180
1.1 Not Detected 3.6 N	2-Propanol	1.1	58	2.8	140
Methylene Chloride 0.46 1.0 1.6 3 Methyl tert-butyl ether 0.23 2.1 0.82 7 trans-1,2-Dichloroethene 0.23 Not Detected 0.91 Not D Hexane 0.23 Not Detected 0.93 Not D 1,1-Dichloroethane 0.23 Not Detected 0.93 Not D 2-Butanone (Methyl Ethyl Ketone) 1.1 4.9 3.4 Not D 2-Butanone (Methyl Ethyl Ketone) 1.1 Not Detected 0.91 Not D 2-Butanone (Methyl Ethyl Ketone) 1.1 Not Detected 0.91 Not D 2-Butanone (Methyl Ethyl Ketone) 1.1 4.9 3.4 Not D Ciss-1,2-Dichloroethene 0.23 Not Detected 0.91 Not D Ciss-1,2-Dichloroethane 0.23 0.067 J 1.1 0.3 Carbon Tetrachloride 0.23 0.092 J 1.4 0.5 Carbon Tetrachloride 0.23 0.092 J 1.4 0.5 Carbon Tetrachloride 0.23	Carbon Disulfide	1.1	0.21 J	3.6	0.65 J
Methyl tert-butyl ether trans-1,2-Dichloroethene 0.23 2.1 0.82 7 dexans-1,2-Dichloroethene 0.23 Not Detected 0.91 Not Detected 1,1-Dichloroethane 0.23 Not Detected 0.93 Not Detected 2-Butanone (Methyl Ethyl Ketone) 1.1 4.9 3.4 3.4 2-Butanone (Methyl Ethyl Ketone) 1.1 4.9 3.4 3.4 5cs-1,2-Dichloroethene 0.23 Not Detected 0.91 Not Detected 6cs-1,2-Dichloroethane 0.23 0.067 J 1.1 0.3 1,1,1-Trichloroethane 0.23 Not Detected 1.2 Not Detected 1,1,1-Trichloroethane 0.23 0.95 J 0.79 3 Carbon Tetrachloride 0.23 0.95 J 0.79 3 Carbon Tetrachloride 0.23 0.092 J 1.4 0.6 2,2,4-Trimethylpentane 1.1 1.8 5.3 8 3-enzene 0.23 4.0 0.73 1 4,2-Dichloroethane 0.23	3-Chloropropene	1.1	Not Detected	3.6	Not Detected
Methyl tert-butyl ether 0.23 2.1 0.82 7 trans-1,2-Dichloroethene 0.23 Not Detected 0.91 Not D Hexane 0.23 1.2 J 0.81 4 1,1-Dichloroethane 0.23 Not Detected 0.93 Not D 2-Butanone (Methyl Ethyl Ketone) 1.1 4.9 3.4 3.4 5cs-1,2-Dichloroethene 0.23 Not Detected 0.91 Not D Cettahydrofuran 1.1 Not Detected 0.91 Not D Chloroform 0.23 0.067 J 1.1 0.3 1,1,1-Trichloroethane 0.23 Not Detected 1.2 Not D Cyclohexane 0.23 0.95 J 0.79 3 Carbon Tetrachloride 0.23 0.092 J 1.4 0.6 2,2,4-Trimethylpentane 1.1 1.8 5.3 8 3enzene 0.23 4.0 0.73 1 4,2-Dichloroethane 0.23 Not Detected 0.93 Not Detected <td>Methylene Chloride</td> <td>0.46</td> <td>1.0</td> <td>1.6</td> <td>3.5</td>	Methylene Chloride	0.46	1.0	1.6	3.5
Hexane	Methyl tert-butyl ether	0.23			7.6
1,1-Dichloroethane 0.23 Not Detected 0.93 Not Detected 0.93 Not Detected 0.94 O.95	rans-1,2-Dichloroethene	0.23	Not Detected	0.91	Not Detected
2-Butanone (Methyl Ethyl Ketone) 1.1 4.9 3.4 63s-1,2-Dichloroethene 0.23 Not Detected 0.91 Not D Cetrahydrofuran 1.1 Not Detected 0.91 1.1 0.3 1,1,1-Trichloroethane 0.23 0.067 J 1.1 0.3 1,1,1-Trichloroethane 0.23 Not Detected 1.2 Not D Cyclohexane 0.23 0.95 J 0.79 3 Carbon Tetrachloride 0.23 0.092 J 1.4 0.5 2,2,4-Trimethylpentane 1.1 1.8 5.3 88 3enzene 0.23 4.0 0.73 1,2-Dichloroethane 0.23 Not Detected 0.93 Not Detected 0.93 Not Detected 0.93 Not D Cyclohexane 0.23 0.77 J 0.94 3 1,2-Dichloroethene 0.23 0.14 J 1.2 0.7 1,2-Dichloropropane 0.23 Not Detected 1.0 Not D Cyclohexane 0.24 Not D Cyclohexane 0.25 Not D Cyclohexane 0.26 Not D Cycloh	Hexane	0.23	1.2 🕽	0.81	4.3
Not Detected Detec	1,1-Dichloroethane	0.23	Not Detected	0.93	Not Detected
cis-1,2-Dichloroethene 0.23 Not Detected 0.91 Not Detected 3.4 Not Detected 3.4 Not Detected 3.4 Not Detected 3.4 Not Detected 1.1 0.3 0.067 J 1.1 0.3 0.07 J 1.1 0.3 0.07 J 1.1 0.3 0.092 J 1.4 0.5 0.093 J 0.0<	2-Butanone (Methyl Ethyl Ketone)	1.1	4.9	3.4	14
Chloroform 0.23 0.067 J 1.1 0.3 1,1,1-Trichloroethane 0.23 Not Detected 1.2 Not Detected Cyclohexane 0.23 0.95 J 0.79 3 Carbon Tetrachloride 0.23 0.092 J 1.4 0.5 Cz,4-Trimethylpentane 1.1 1.8 5.3 8 Senzene 0.23 4.0 0.73 6 3enzene 0.23 4.0 0.73 6 1,2-Dichloroethane 0.23 Not Detected 0.93 Not Detected 1,2-Dichloropropane 0.23 0.14 J 1.2 0.7 1,2-Dichloropropane 0.23 0.17 J 0.82 0.6 3cromodichloromethane 0.23 0.17 J 0.82 0.6 3cromodichloropropene 0.23 0.65 0.94 2 4-Methyl-2-pentanone 0.23 0.65 0.94 2 Foluene 0.23 Not Detected 1.0 Not Detected 1,1,2-Trichloroethane<		0.23	Not Detected		Not Detected
1,1,1-Trichloroethane	Tetrahydrofuran	1.1	Not Detected J	3.4	Not Detected
Cyclohexane 0.23 0.95 J 0.79 3 Carbon Tetrachloride 0.23 0.092 J 1.4 0.5 2,2,4-Trimethylpentane 1.1 1.8 5.3 8 Benzene 0.23 4.0 0.73 1 1,2-Dichloroethane 0.23 Not Detected 0.93 Not Detected 1,2-Dichloropropane 0.23 0.77 J 0.94 3 1,2-Dichloropropane 0.23 0.14 J 1.2 0.7 1,2-Dichloropropane 0.23 0.17 J 0.82 0.6 1,4-Dioxane 0.23 0.17 J 0.82 0.6 3romodichloromethane 0.23 0.17 J 0.82 0.6 4-Methyl-2-pentanone 0.23 0.65 0.94 2 Foluene 0.23 Not Detected 1.0 Not Detected 1,1,2-Trichloroethane 0.23 Not Detected 1.6 Not Detected 1,1,2-Trichloroethane 0.23 Not Detected 1.6 Not Detected	Chloroform	0.23	0.067 J	1.1	0.32 J
Carbon Tetrachloride 0.23 0.092 J 1.4 0.5 2,2,4-Trimethylpentane 1.1 1.8 5.3 8 Benzene 0.23 4.0 0.73 6 1,2-Dichloroethane 0.23 Not Detected 0.93 Not Detected 1,2-Dichloropropane 0.23 0.14 J 1.2 0.7 1,2-Dichloropropane 0.23 Not Detected 1.0 Not Detected 1,4-Dioxane 0.23 0.17 J 0.82 0.6 3-romodichloromethane 0.23 Not Detected 1.5 Not Detected 4-Methyl-2-pentanone 0.23 0.65 0.94 2 7-oluene 0.23 Not Detected 1.0 Not Detected 1,1,2-Trichloropropene 0.23 Not Detected 1.0 Not Detected 1,1,2-Trichloroethane 0.23 Not Detected 1.6 Not Detected	1,1,1-Trichloroethane	0.23	Not Detected	1.2	Not Detected
Carbon Tetrachloride 0.23 0.092 J 1.4 0.5 2,2,4-Trimethylpentane 1.1 1.8 5.3 8 Benzene 0.23 4.0 0.73 6 1,2-Dichloroethane 0.23 Not Detected 0.93 Not Detected 1,2-Dichloropropane 0.23 0.14 J 1.2 0.7 1,2-Dichloropropane 0.23 Not Detected 1.0 Not Detected 1,4-Dioxane 0.23 0.17 J 0.82 0.6 3-romodichloromethane 0.23 Not Detected 1.5 Not Detected 4-Methyl-2-pentanone 0.23 0.65 0.94 2 7-oluene 0.23 Not Detected 1.0 Not Detected 1,1,2-Trichloropropene 0.23 Not Detected 1.0 Not Detected 1,1,2-Trichloroethane 0.23 Not Detected 1.6 Not Detected	Cyclohexane	0.23	0.95 7	0.79	3.3
2,2,4-Trimethylpentane 1.1 1.8 5.3 8 Benzene 0.23 4.0 0.73 1 1,2-Dichloroethane 0.23 Not Detected 0.93 Not Detected 1,2-Dichloroethene 0.23 0.14 J 1.2 0.7 1,2-Dichloropropane 0.23 Not Detected 1.0 Not Detected 1,4-Dioxane 0.23 0.17 J 0.82 0.6 3-Gromodichloromethane 0.23 Not Detected 1.5 Not Detected 1-Methyl-2-pentanone 0.23 0.65 0.94 2 1-Methyl-2-pentanone 0.23 0.65 0.94 2 1-Methyl-2-pentanone 0.23 Not Detected 1.0 Not Detected 1,1,2-Trichloroethane 0.23 Not Detected 1.0 Not Detected 1,1,2-Trichloroethane 0.23 Not Detected 1.6 Not Detected	Carbon Tetrachloride	0.23	•		0.58 J
Benzene 0.23 4.0 0.73 1 1,2-Dichloroethane 0.23 Not Detected 0.93 Not Detected 0.94 3 3 3 3 3 3 3 3 3	2,2,4-Trimethylpentane	1.1	1.8		8.3
1,2-Dichloroethane	Benzene	0.23	4.0	0.73	13
Heptane	1,2-Dichloroethane	0.23	Not Detected		Not Detected
Trichloroethene 0.23 0.14 J 1.2 0.7 1,2-Dichloropropane 0.23 Not Detected 1.0 Not Detected 1,4-Dioxane 0.23 0.17 J 0.82 0.6 3romodichloromethane 0.23 Not Detected 1.5 Not Detected 4-Methyl-2-pentanone 0.23 0.65 0.94 2 Foluene 0.23 Not Detected 1.0 Not Detected 1,1,2-Trichloropropene 0.23 Not Detected 1.0 Not Detected 1,1,2-Trichloroethane 0.23 Not Detected 1.6 Not Detected	Heptane	0.23	0.77 ゴ	0.94	3.1
1,2-Dichloropropane 0.23 Not Detected 1.0 Not Detected 1,4-Dioxane 0.23 0.17 J 0.82 0.6 3romodichloromethane 0.23 Not Detected 1.5 Not Detected 1.5 Not Detected 1.0 Not Detected 1.0 Not Detected 1.0 Not Detected 0.86 2 1-Methyl-2-pentanone 0.23 0.65 0.94 2 2 1-foluene 0.23 Not Detected 1.0 Not Detected 1.0 Not Detected 1.0 Not Detected 1.2 Not Detected 1.2 Not Detected 1.6 Not Detected 1.0 Not Detected 1.6 Not Detected 1.6 Not Detected 1.6 Not Detected 1.0	•				0.73 J
1,4-Dioxane 0.23 0.17 J 0.82 0.6 Bromodichloromethane 0.23 No Detected 1.5 Not Detected 4-Methyl-2-pentanone 0.23 0.65 0.94 2 Foluene 0.23 0.65 0.86 2 rans-1,3-Dichloropropene 0.23 Not Detected 1.0 Not Detected 1,1,2-Trichloroethane 0.23 Not Detected 1.6 Not Detected	1,2-Dichloropropane	0.23			Not Detected
Stromodichloromethane 0.23 Not Detected 1.5 Not Detected 1.5 Not Detected 1.6 Not Detected 1.7 Not Detected 1.7 Not Detected 1.8 Not Detected 1.9	1,4-Dioxane				0.62 J
1.0 Not December 1.1,1,2-Trichloroethane 1.2 Not December 1.2 Not December 1.3 Not December 1.4 Not December 1.5 Not December 1.6 Not December 1.7 Not December 1.	Bromodichloromethane				Not Detected
1-Methyl-2-pentanone	cis-1,3-Dichloropropene	0.23	OF NOT DONCE TO		Not Detected
Toluene 0.23 Sinte Difference 0.86 2 rans-1,3-Dichloropropene 0.23 Not Defeated 1.0 Not Defeated 1.1.2 Not Defeated 1.2 Not Defeated 1.2 Not Defeated 1.3 Not Defeated 1.4 Not Defeated 1.5 Not Defeated 1.6 Not	• •		0.65		2.7
rans-1,3-Dichloropropene 0.23 Not Enterted 1.0 Not De 1,1,2-Trichloroethane 0.23 Not DetEcted 1.2 Not De 1.2 N		0.23	Salafaeblicae 18		25
Tetrachloroethene 0.23 Not Detected 1.2 Not Detected 1.6		0.23	Not Reflected	1	Not Detected
Tetrachloroethene 0.23 Not Detected 1.6 Not Detected 1.6		0.23	NocDetians		Not Detected
				/	Not Detected
chickenione i.i who Determine 4.7 Not Determine 4.7	2-Hexanone	1.1	Mor Detected 1	4.7	Not Detected



Client Sample ID: B6-3SS Lab ID#: 1502113AR1-02A

MODIFIED EPA METHOD TO-15 GC/MS FULL SCAN

File Name: Dil. Factor:	v021016r1 2.29		te of Collection: 2/4/ te of Analysis: 2/10/1	
Compound	Rpt. Limit (ppbv)	Amount (ydag)	Rpt. Limit (ug/m3)	Amount (ug/m3)

Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Dibromochloromethane	0.23	Not Detected	2.0	Not Detected
1,2-Dibromoethane (EDB)	0.23	Not Detected	1.8	Not Detected
Chlorobenzene	0.23	0.11 J	1.0	0.50 J
Ethyl Benzene	0.23	12	0.99	54
m,p-Xylene	0.23	1.4	0.99	5.9
o-Xylene	0.23	0.42	0.99	1.8
Styrene	0.23	11	0.98	45
Bromoform	0.23	Not Detected	2.4	Not Detected
Cumene	0.23	0.10 J	1.1	0.51 J
1,1,2,2-Tetrachloroethane	0.23	Not Detected	1.6	Not Detected
Propylbenzene	0.23	0.16 J J	1.1	0.79 J
4-Ethyltoluene	0.23	0.22 J	1.1	1.0 J
1,3,5-Trimethylbenzene	0.23	Not Detected	1.1	Not Detected
1,2,4-Trimethylbenzene	0.23	0.27	1.1	1.4
1,3-Dichlorobenzene	0.23	Not Detected	1.4	Not Detected
1,4-Dichlorobenzene	0.23	Not Detected	1.4	Not Detected
alpha-Chlorotoluene	1.1	Not Detected	5.9	Not Detected
1,2-Dichlorobenzene	0.23	Not Detected	1.4	Not Detected
1,2,4-Trichlorobenzene	1.1	Not Detected	8.5	Not Detected
Hexachlorobutadiene	1.1	Not Detected UJ 3	12	Not Detected UJ
Naphthalene	1.1	Not Detected	6.0	Not Detected

J = Estimated value.

UJ = Analyte associated with low bias in the CCV and/or LCS. Container Type: 1 Liter Summa Canister (100% Certified)

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	102	70-130
Toluene-d8	98	70-130
4-Bromofluorobenzene	90	70-130



J0 = Estimated value due to bias in the CCV.



Client Sample ID: B6-3SSD Lab ID#: 1502113AR1-03A

MODIFIED EPA METHOD TO-15 GC/MS FULL SCAN

File Name: Dil. Factor:	v021018r1 2.29		of Collection: 2/4 of Analysis: 2/10/	
	Rpt. Limit	Amount	Rpt. Limit	Amount
Compound	(ppbv)	(ppbv)	(ug/m3)	(ug/m3)
Freon 12	0.23	0.47	1.1	2.3
Freon 114	0.23	Not Detected	1.6	Not Detected
Chloromethane	1.1	0.74 J	2.4	1.5 J
Vinyl Chloride	0.23	Not Detected	0.58	Not Detected
1,3-Butadiene	0.23	Not Detected	0.51	Not Detected
Bromomethane	1.1	Not Detected	4.4	Not Detected
Chloroethane	1.1	Not Detected	3.0	Not Detected
Freon 11	0.23	0.21 J	1.3	1.2 J
Ethanol	1.1	16 JO J	2.2	30 J0
Freon 113	0.23	0.059 J	1.8	0.45 J
1,1-Dichloroethene	0.23	Not Detected	0.91	Not Detected
Acetone	1.1	76 1	2.7	180
2-Propanol	1,1	61	2.8	150
Carbon Disulfide	1.1	Not Detected	3.6	Not Detected
3-Chloropropene	1.1	Not Detected	3.6	Not Detected
Methylene Chloride	0.46	1.0	1.6	3.6
Methyl tert-butyl ether	0.23	2.4	0.82	8.7
rans-1,2-Dichloroethene	0.23	0.063 J	0.91	0.25 J
-lexane	0.23	0.51 🎝	0.81	1.8
1,1-Dichloroethane	0.23	Not Detected	0.93	Not Detected
2-Butanone (Methyl Ethyl Ketone)	1.1	4.8	3.4	14
cis-1,2-Dichloroethene	0.23	Not Detected	0.91	Not Detected
Tetrahydrofuran	1.1	0.26 J ゴ	3.4	0.75 J
Chloroform	0.23	0.050 J	1.1	0.24 J
1,1,1-Trichloroethane	0.23	Not Detected	1.2	Not Detected
Cyclohexane	0.23	0.38 ブ	0.79	1.3
Carbon Tetrachloride	0.23	0.078 J	1.4	0.49 J
2,2,4-Trimethylpentane	1.1	Not Detected	5.3	Not Detected
Benzene	0.23	1.2	0.73	3.7
1,2-Dichloroethane	0.23	Not Detected	0.93	Not Detected
- Heptane	0.23	0.28 🤰	0.94	1.1
Trichloroethene	0.23	0.16 J	1.2	0.89 J
1,2-Dichloropropane	0.23	Not Detected	1.0	Not Detected
1,4-Dioxane	0.23	0.22 J	0.82	0.80 J
Bromodichloromethane	0.23	Not Detected	1.5	Not Detected
cis-1,3-Dichloropropene	0.23	Not Detected	1.0	Not Detected
4-Methyl-2-pentanone	and the second s	0.68	0.94	2.8
Foluene	SOCIADO SE	7.0	0.86	26
rans-1,3-Dichloropropene /		Not Detected	1.0	Not Detected
1,1,2-Trichloroethane	fael Infah@	Not Detected	1.2	Not Detected
Tetrachloroethene	Ménde 20.23	Not Detected	1.6	
Tetrachloroethene 2-Hexanone	IC # 18481	0.35 J		Not Detected
2-mexamone	IC # 10=491	U.33 J ✓	4.7	1.4 J

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Client Sample ID: B6-3SSD Lab ID#: 1502113AR1-03A

MODIFIED EPA METHOD TO-15 GC/MS FULL SCAN

File Name: Dil. Factor:	v021018r1 2.29	Date of Collection: 2/4/15 4:37:00 PM Date of Analysis: 2/10/15 08:21 PM		
Compound	Rpt. Limit (ppby)	Amount	Rpt. Limit	Amount
Dibromochloromethane	0.53 (bbps)	(ppbv)	(ug/m3)	(ug/m3)

Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Dibromochloromethane	0.23	Not Detected	2.0	Not Detected
1,2-Dibromoethane (EDB)	0.23	Not Detected	1.8	Not Detected
Chlorobenzene	0.23	0.12 J	1.0	0.55 J
Ethyl Benzene	0.23	13	0.99	57
m,p-Xylene	0.23	0.98	0.99	4.2
o-Xylene	0.23	0.43	0.99	1.9
Styrene	0.23	12	0.98	49
Bromoform	0.23	Not Detected	2.4	Not Detected
Cumene	0.23	0.060 J	1.1	0.29 J
1,1,2,2-Tetrachloroethane	0.23	Not Detected	1.6	Not Detected
Propylbenzene	0.23	0.057 J 🤈	1.1	0.28 J
4-Ethyltoluene	0.23	0.21 J	1.1	1.0 J
1,3,5-Trimethylbenzene	0.23	Not Detected	1.1	Not Detected
1,2,4-Trimethylbenzene	0.23	0.24	1.1	1.2
1,3-Dichlorobenzene	0.23	Not Detected	1.4	Not Detected
1,4-Dichlorobenzene	0.23	Not Detected	1.4	Not Detected
alpha-Chlorotoluene	1.1	Not Detected	5.9	Not Detected
1,2-Dichlorobenzene	0.23	Not Detected	1.4	Not Detected
1,2,4-Trichlorobenzene	1.1	Not Detected	8.5	Not Detected
Hexachlorobutadiene	1.1	Not Detected UJ 3	12	Not Detected UJ
Naphthalene	1.1	Not Detected	6.0	Not Detected

J = Estimated value.

UJ = Analyte associated with low bias in the CCV and/or LCS. Container Type: 1 Liter Summa Canister (100% Certified)

Surrogates	%Recovery	Method Limits	
1,2-Dichloroethane-d4	104	70-130	
Toluene-d8	97	70-130	
4-Bromofluorobenzene	90	70-130	



J0 = Estimated value due to bias in the CCV.



Client Sample ID: B6-4SS Lab ID#: 1502113AR1-04A

MODIFIED EPA METHOD TO-15 GC/MS FULL SCAN

File Name: Dil. Factor:	v021019r1 12.4	Date of Collection: 2/6/15 2:22:00 PM Date of Analysis: 2/10/15 08:55 PM		
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Freon 12	1.2	Not Detected	6.1	Not Detected
Freon 114	1.2	Not Detected	8.7	Not Detected
Chloromethane	6.2	1.1 J	13	2.2 J
Vinyl Chloride	1.2	Not Detected	3.2	Not Detected
1,3-Butadiene	1.2	Not Detected	2.7	Not Detected
Bromomethane	6.2	Not Detected	24	Not Detected
Chloroethane	6.2	Not Detected	16	Not Detected
Freon 11	1.2	Not Detected	7.0	Not Detected
Ethanol	6.2	89 JO J	12	170 J0
Freon 113	1.2	Not Detected	9.5	Not Detected
1,1-Dichloroethene	1.2	Not Detected	4.9	Not Detected
Acetone	6.2	170 J	15	400
2-Propanol	6.2	1900 E	15	4700 E
Carbon Disulfide	6.2	Not Detected	19	Not Detected
3-Chloropropene	6.2	Not Detected	19	Not Detected
Methylene Chloride	2.5	1.1 J	8.6	3.7 J
Methyl tert-butyl ether	1.2	Not Detected	4.5	Not Detected
trans-1,2-Dichloroethene	1.2	Not Detected	4.9	Not Detected
Hexane	1.2	1.1 J	4.4	4.0 J
1,1-Dichloroethane	1.2	Not Detected	5.0	Not Detected
2-Butanone (Methyl Ethyl Ketone)	6.2	1.6 J	18	4.7 J
cis-1,2-Dichloroethene	1.2	Not Detected	4.9	Not Detected
Tetrahydrofuran	6.2	Not Detected	18	Not Detected
Chloroform	1.2	0.25 J	6.0	1.2 J
1,1,1-Trichloroethane	1.2	Not Detected	6.8	Not Detected
Cyclohexane	1.2	Not Detected	4.3	Not Detected
Carbon Tetrachloride	1.2	Not Detected	7.8	Not Detected
2,2,4-Trimethylpentane	6.2	Not Detected	29	Not Detected
Benzene	1.2	Not Detected	4.0	Not Detected
1,2-Dichloroethane	1.2	Not Detected	5.0	Not Detected
Heptane	1.2	Not Detected	5.1	Not Detected
Trichloroethene	1.2	Not Detected	6.7	Not Detected
1,2-Dichloropropane	1.2	Not Detected	5.7	Not Detected
1,4-Dioxane	1.2	Not Detected	4.5	Not Detected
Bromodichloromethane	1.2	Not Detected	8.3	Not Detected
cis-1,3-Dichloropropene	1.2	Not Detected	5.6	Not Detected
4-Methyl-2-pentanone		0.43 J	5.1	1.8 J
Toluene		2.5	4.7	9.5
trans-1,3-Dichloropropene	fael Infinite	Not Detected	5.6	Not Detected
	Méndez 18	Not Detected	6.8	Not Detected
Tetrachloroethene \\ \tilde{\chi} \\ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \	C # 1885	Not Detected	8.4	Not Detected
2-Hexanone	/&/	Not Detected	25	Not Detected

Page 1



Client Sample ID: B6-4SS Lab ID#: 1502113AR1-04A

MODIFIED EPA METHOD TO-15 GC/MS FULL SCAN

File Name: Dil. Factor:	v021019r1 12.4		Pate of Collection: 2/6/15 2:22:00 PM Pate of Analysis: 2/10/15 08:55 PM		
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)	
Dibromochloromethane	1.2	Not Detected	10	Not Detected	
1,2-Dibromoethane (EDB)	1.2	Not Detected	9.5	Not Detected	
Chlorobenzene	1.2	Not Detected	5.7	Not Detected	
Ethyl Benzene	1.2	1.8	5.4	8.0	
m,p-Xylene	1.2	11	5.4	48	
o-Xylene	1.2	4.1	5.4	18	
Styrene	1.2	Not Detected	5.3	Not Detected	
Bromoform	1.2	Not Detected	13	Not Detected	
Cumene	1.2	- 0.30 J	6.1	1.5 J	
1,1,2,2-Tetrachloroethane	1.2	Not Detected	8.5	Not Detected	
Propylbenzene	1.2	0.79 J	6.1	3.9 J	
4-Ethyltoluene	1.2	2.4	6.1	12	
1,3,5-Trimethylbenzene	1.2	1.1 J	6.1	5.3 J	
1,2,4-Trimethylbenzene	1.2	3.7	6.1	18	
1,3-Dichlorobenzene	1.2	Not Detected	7.4	Not Detected	
1,4-Dichlorobenzene	1.2	Not Detected	7.4	Not Detected	
alpha-Chiorotoluene	6.2	Not Detected	32	Not Detected	
1,2-Dichlorobenzene	1.2	Not Detected	7.4	Not Detected	
1,2,4-Trichlorobenzene	6.2	Not Detected	46	Not Detected	
Hexachlorobutadiene	6.2	Not Detected UJ 3	66	Not Detected UJ	

J = Estimated value.

Naphthalene

UJ = Analyte associated with low bias in the CCV and/or LCS. Container Type: 1 Liter Summa Canister (100% Certified)

		method
Surrogates	%Recovery	Limits
1,2-Dichloroethane-d4	97	70-130
Toluene-d8	97	70-130
4-Bromofluorobenzene	88	70-130

Not Detected

32

6.2

Not Detected

J0 = Estimated value due to bias in the CCV.

E = Exceeds instrument calibration range.

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024 BG-355 37729 2/4/15 1632	5		
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OHA B6-455 11427 2/6/15 1422 30	5		
	 		
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	Project Number:1502113AR1
	Date:02/04-06/2015
REVIEW OF VOLATILE ORGATHE ORGATHE following guidelines for evaluating volatile organics wactions. This document will assist the reviewer in using prodecision and in better serving the needs of the data users. The USEPA data validation guidance documents in the follow "Compendium Method TO-15. Determination of Volatile Org Specially-Prepared Canisters and Analyzed By Gas Chellow January, 1999"; USEPA Hazardous Waste Support Brand Analysis of Ambient Air in Canisters by Method TO-15, (SOF QC criteria and data validation actions listed on the data review document, unless otherwise noted. The hardcopied (laboratory name) _EurofinsAir_Toxics	ANIC PACKAGE vere created to delineate required validation ofessional judgment to make more informed the sample results were assessed according to bring order of precedence: QC criteria from ganic Compounds (VOCs) In Air Collected In romatography/Mass Spectrometry (GC/MS), th. Validating Air Samples. Volatile Organic P # HW-31. Revision #4. October, 2006). The ew worksheets are from the primary guidance
reviewed and the quality control and performance data summa	arized. The data review for VOCs included:
Lab. Project/SDG No.:1502113AR1 No. of Samples:4	Sample matrix:Air
Trip blank No.: Field blank No.: Equipment blank No.: Field duplicate No.:B6-3SS/B6-3SSD	
X Data CompletenessX Holding TimesX GC/MS TuningX Internal Standard PerformanceX BlanksX Surrogate RecoveriesN/A_ Matrix Spike/Matrix Spike Duplicate	X Laboratory Control SpikesX Field DuplicatesX CalibrationsX Compound IdentificationsX Compound QuantitationX Quantitation Limits
Overall Comments:_VOCs_by_method_TO-15	
Definition of Qualifiers: J- Estimated results U- Compound not detected R- Rejected data UJ- Estimated nondetect Reviewer: Date: 06/30/2015	

DATA REVIEW WORKSHEETS

DATA COMPLETENESS

MISSING INFORMATION	DATE LAB. CONTACTED	DATE RECEIVED
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All criteria were metX
Criteria were not met
and/or see below

HOLDING TIMES

The objective of this parameter is to ascertain the validity of the results based on the holding time of the sample from time of collection to the time of analysis.

Complete table for all samples and note the analysis and/or preservation not within criteria

SAMPLE ID	DATE SAMPLED	DATE ANALYZED	pН	ACTION
,	All samples analyzed w	ithin the recommended	l method	holding time

Criteria

Aqueous samples – 14 days from sample collection for preserved samples (pH \leq 2, 4°C), no air bubbles.

Aqueous samples – 7 days from sample collection for unpreserved samples, 4°C, no air bubbles. Soil samples- 7 days from sample collection.

Cooler temperature (Criteria: 4 ± 2 °C): N/A – summa canisters

Actions

If the VOCs vial(s) have air bubbles, estimate positive results (J) and reject nondetects (R).

If the % solids of soil samples is 10-50%, estimates positive results (J) and nondetects (UJ)

If the % solid of soil samples is < 10%, estimate positive results (J) and reject nondetects (R).

If holding times are exceeded but < 14 days beyond criteria, estimate positive results (J) and nondetects (UJ).

If holding times are exceeded but < 28 days beyond criteria, estimate positive results (J) and reject nondetects (R).

If holding times are grossly exceeded (> 28 days beyond criteria), reject all results (R).

If samples were not iced or if the ice were melted (> 10°C), estimate positive results (J) and nondetects (UJ).

DATA REVIEW WORKSHEETS

		Crit	All criteria were metX teria were not met see below
GC/MS TUNING			
The assessment o standard tuning QC		o determine if the sample instru	mentation is within the
_XThe BFB p	erformance results were	e reviewed and found to be within	the specified criteria.
_XBFB tuning	was performed for ever	ry 24 hours of sample analysis.	
f no, use professi qualified or rejected		mine whether the associated dat	ta should be accepted,
List	the	samples	affected:
	Va.		

If mass calibration is in error, all associated data are rejected.

All criteria were metX_	_
Criteria were not met	
and/or see below	

CALIBRATION VERIFICATION

Compliance requirements for satisfactory instrument calibration are established to ensure that the instrument is capable of producing and maintaining acceptable quantitative data.

Date of initial calibration:	_01/23/15
Dates of continuing calibration:	02/10/15
Instrument ID numbers:MS	SD-V
Matrix/Level:/	\ir/low

DATE	LAB FILE	CRITERIA OUT RFs, %RSD, %D, r	COMPOUND	SAMPLES AFFECTED
Initial and	continuing calib	ration met the method pe	 rformance criteria except f	or the followings:
02/10/15	v021002.d	- 39 % RSD	Acetone	All samples.
		39 % D	Ethanol	•
		-32 % D	Hexachlorobutadiene	1.

Criteria

All RFs must be > 0.05 regardless of method requirements for SPCC.

All %RSD must be \leq 15 % regardless of method requirements for CCC.

All %Ds must be \leq 30% regardless of method requirements for CCC.

Method TO-15 does not specify criterion for the curve correlation coefficient (r). A limit for r of \geq 0.995 has therefore been utilized as professional judgment.

Actions

If any compound has an initial RF or a continuing RF of < 0.05, estimate positive results (J) and reject nondetects (R), regardless of method requirements.

If any compound has a %RSD > 15%, estimate positive results (J) and use professional judgment to qualify nondetects.

If any compound has a %RSD > 90%, estimate positive results (J) and reject nondetects (R).

If any compound has a % D > 30%, estimate positive results (J) and reject nondetects (R).

If any compound has a % D > 30%, estimate positive results (J) and nondetects (UJ).

If any compound has a % D > 90%, estimate positive results (J) and reject nondetects (R).

If any compound has r < 0.995, estimate positive results and nondetects.

A separate worksheet should be filled for each initial curve

All criteria were met
Criteria were not met
and/or see belowX

V A. BLANK ANALYSIS RESULTS (Sections 1 & 2)

The assessment of the blank analysis results is to determine the existence and magnitude of contamination problems. The criteria for evaluation of blanks apply only to blanks associated with the samples, including trip, equipment, and laboratory blanks. If problems with any blanks exist, all data associated with the case must be carefully evaluated to determine whether or not there is an inherent variability in the data for the case, or if the problem is an isolated occurrence not affecting other data.

List the contamination in the blanks below. High and low levels blanks must be treated separately.

Laboratory blanks

DATE ANALYZED	LAB ID	LEVEL/ MATRIX	COMPOUND	CONCENTRATION UNITS
All_metho	d_blank_meeth_me			he_followings:
_02/10/15	1502113AR1-05A	Air/low	Acetone	0.32_ppbv
Field/Equipmen		***************************************		
DATE ANALYZED	LAB ID	LEVEL/ MATRIX	COMPOUND	CONCENTRATION UNITS
No_field/trip/eq	uipment_blanks_ar	nalyzed_with	n_this_data_package	

All criteria were metX
Criteria were not met
and/or see below

VB. BLANK ANALYSIS RESULTS (Section 3)

Blank Actions

Action Levels (ALs) should be based upon the highest concentration of contaminant determined in any blank. Do not qualify any blank with another blank. The ALs for samples which have been diluted should be corrected for the sample dilution factor and/or % moisture, where applicable. No positive sample results should be reported unless the concentration of the compound in the samples exceeds the ALs:

ALs = 10x the amount of common contaminants (methylene chloride, acetone, 2-butanone, and toluene)

ALs = 5x for any other compounds

Specific actions are as follows:

If the concentration is < sample quantitation limit (SQL) and \le AL, report the compound as not detected (U) at the SQL.

If the concentration is \geq SQL but \leq AL, report the compound as not detected (U) at the reported concentration.

If the concentration is \geq SQL and > AL, report the concentration unqualified.

Notes:

High and low level blanks must be treated separately

Compounds qualified "U" for blank contamination are still considered "hits" when qualifying for calibration criteria.

CONTAMINATION SOURCE/LEVEL	COMPOUND	CONC/UNITS	AL/UNITS	SQL	AFFECTED SAMPLES
					,

All criteria were metX	
Criteria were not met	
and/or see below	

SURROGATE SPIKE RECOVERIES

d4

Laboratory performance of individual samples is established by evaluation of surrogate spike recoveries. All samples are spiked with surrogate compounds prior to sample analysis. The accuracy of the analysis is measured by the surrogate percent recovery. Since the effects of the sample matrix are frequently outside the control of the laboratory and may present relatively unique problems, the validation of data is frequently subjective and demands analytical experience and professional judgment.

List the percent recoveries (%Rs) which do not meet the criteria for surrogate recovery.

Matrix: solid/aqueous

SAMPLE ID

SURROGATE COMPOUND

ACTION

1,2-DICHLOROETHANE-

Toluene- 4-BFB

d8

Surrogate_rec	overies_w	thin_labora	tory_contro	l_limits		
						
-						
C Limits* (Air)						
LL_to_L	JL 70	to_130_		70 to 130	70 to 130	

- * QC limits are laboratory in-house performance criteria, LL = lower limit, UL = upper limit.
- * If QC limits are not available, use limits of 80 120 % for aqueous and 70 130 % for solid samples.

Actions:

QUALITY	%R < 10%	%R = 10% - LL	%R > UL
Positive results	J	J	J
Nondetects results	R	UJ	Accept

Surrogate action should be applied:

If one or more surrogate in the VOC fraction is out of specification, but has a recovery of > 10%.

If any one surrogate in a fraction shows < 10 % recovery.

All criteria were met
Criteria were not met
and/or see belowN/A

VII. B MATRIX SPIKE/MATRIX SPIKE DUPLICATE

MS/MSD - Unspiked Compounds

It should be noted that Method TO-15 does not specify a MS/MSD criteria for the unspiked compounds in the sample. A %RSD of < 50% has therefore been utilized as professional judgment.

If all target analytes were spiked in the MS/MSD, this review element is not applicable.

List the %RSD of the compounds which do not meet the criteria.

Sample ID:			Matrix/Le	vel/Unit:	
COMPOUND	SAMPLE CONC.	MS CONC.	MSD CONC.	% RSD	ACTION
A-2					
					-
	£				

Actions:

^{*} If the % RSD > 50, qualify the positive result in the unspiked samples as estimated (J).

^{*} If the % RSD is not calculated (NC) due to nondetected value, use professional judgment to qualify the data.

All criteria were metX
Criteria were not met
and/or see below

VIII. LABORATORY CONTROL SAMPLE (LCS) ANALYSIS

This data is generated to determine accuracy of the analytical method for various matrices.

1. LCS Recoveries Criteria

Where LCS spiked with the same analyte at the same concentrations as the MS/MSD? Yes or No. If no make note in data review memo.

List the %R of compounds which do not meet the criteria

	LCS ID	COMPOUND	% R	QC LIMIT
LCS/L0	CSD_%_recoverie	s_and_RPD_within_labora	atory_control_limits	

		101		
			MA	

- * QC limits are laboratory in-house performance criteria, LL = lower limit, UL = upper limit.
- * If QC limits are not available, use limits of 70 130 %.

Actions:

QUALITY	%R < LL	%R > UL
Positive results	J	J
Nondetects results	R	Accept

All analytes in the associated sample results are qualified for the following criteria.

If 25 % of the LCS recoveries were < LL (or 70 %), qualify all positive results (j) and reject nondetects (R).

If two or more LCS were below 10 %, qualify all positive results as (J) and reject nondetects (R).

2. Frequency Criteria:

Where LCS analyzed at the required frequency and for each matrix? Yes or No.

If no, the data may be affected. Use professional judgment to determine the severity of the effect and qualify data accordingly. Discuss any actions below and list the samples affected.

		All criteria were metX Criteria were not met and/or see below
IX.	LABORATORY DUPLICATE PRECISION	
	Sample IDs:LCS/LCSD	Matrix:Air

Laboratory duplicates samples may be taken and analyzed as an indication of overall precision. These analyses measure both field and lab precision; therefore, the results may have more variability than laboratory duplicates which only laboratory performance. It is also expected that soil duplicate results will have a greater variance than water matrices due to difficulties associated with collecting identical field duplicate samples.

The project QAPP should be reviewed for project-specific information.

Suggested criteria: RPD <u>+</u> 25% for air samples. If both samples and duplicate are <5 SQL, the RPD criteria is doubled.

COMPOUND	SQL	SAMPLE CONC.	DUPLICATE CONC.	RPD	ACTION	
						-
N N	RPD Within lab	oratory and (generally accept	able con	trol limits.	· · · · · · · · · · · · · · · · · · ·

Actions:

Qualify as estimated positive results (J) and nondetects (UJ) for the compound that exceeded the above criteria. For organics, only the sample and duplicate will be qualified.

If an RPD cannot be calculated because one or both of the sample results is not detected, the following actions apply:

If one sample result is not detected and the other is greater than 5x the SQL qualify (J/UJ).

If one sample value is not detected and the other is greater than 5x the SQL and the SQLs for the sample and duplicate are significantly different, use professional judgment to determine if qualification is appropriate.

If one sample value is not detected and the other is less than 5x, use professional judgment to determine if qualification is appropriate.

If both sample and duplicate results are not detected, no action is needed.

All criteria were met
Criteria were not met
and/or see belowX

IX. FIELD DUPLICATE PRECISION

Sample IDs:	B6-3SS/B6-3SSD	Matrix:Air
-------------	----------------	------------

Field duplicates samples may be taken and analyzed as an indication of overall precision. These analyses measure both field and lab precision; therefore, the results may have more variability than laboratory duplicates which only laboratory performance. It is also expected that soil duplicate results will have a greater variance than water matrices due to difficulties associated with collecting identical field duplicate samples.

The project QAPP should be reviewed for project-specific information.

Suggested criteria: RPD \pm 25% for air samples. If both samples and duplicate are <5 SQL, the RPD criteria is doubled.

COMPOUND	SQL	SAMPLE CONC.	DUPLICATE CONC.	RPD	ACTION
Hexane	0.03452	1.2	0.51	81 %	Qualify results (J) in sample and duplicate
Tetrahydrofuran	-	ND	0.26	NR	Qualify results (J) in sample and duplicate
Cyclohexane	0.03469	0.95	0.38	86 %	Qualify results (J) in sample and duplicate
Propylbenzene	0.02205	0.16	0.057	95 %	Qualify results (J) in sample and duplicate
Heptane	0.02709	0.77	0.28	96 %	Qualify results (J) in sample and duplicate
2-Hexanone	-	ND	0.35	93 %	Qualify results (J) in sample and duplicate

Actions:

Qualify as estimated positive results (J) and nondetects (UJ) for the compound that exceeded the above criteria. For organics, only the sample and duplicate will be qualified.

If an RPD cannot be calculated because one or both of the sample results is not detected, the following actions apply:

If one sample result is not detected and the other is greater than 5x the SQL qualify (J/UJ).

If one sample value is not detected and the other is greater than 5x the SQL and the SQLs for the sample and duplicate are significantly different, use professional judgment to determine if qualification is appropriate.

If one sample value is not detected and the other is less than 5x, use professional judgment to determine if qualification is appropriate.

If both sample and duplicate results are not detected, no action is needed.

All criteria were metX
Criteria were not met
and/or see below

X. INTERNAL STANDARD PERFORMANCE

The assessment of the internal standard (IS) parameter is used to assist the data reviewer in determining the condition of the analytical instrumentation.

List the internal standard area of samples which do not meet the criteria.

- * Area of +40% or -40% of the IS area in the associated calibration standard.
- * Retention time (RT) within \pm 0.06 seconds of the IS area in the associated calibration standard.

DATE	SAMPLE ID	IS OUT	IS AREA	ACCEPTABLE RANGE	ACTION
	tandard_area_and_reration_standards	etention_times_	within_laboratory	_control_limits_for_	_both_samples
***		· .			
Actions:					

1. IS actions should be applied to the compound quantitated with the out-of-control ISs

QUALITY	IS AREA < -40%	IS AREA > + 40%
Positive results	J	J
Nondetected results	R	ACCEPT

2. If a IS retention time varies more than 0.330 seconds, the chromatographic profile for that sample must be examined to determine if any false positive or negative exists. For shifts of a large magnitude, the reviewer may consider partial or total rejection of the data for the sample fraction.

All criteria were metX
Criteria were not met
and/or see below

XII. SAMPLE QUANTITATION

The sample quantitation evaluation is to verify laboratory quantitation results. In the space below, please show a minimum of one sample calculation:

1502113AR1-01A

MTBE RF = 4.45779

[] = (1466385)(5.0)/(190063)(4.45779)

= 8.65 ppbv OK

All criteria were met
Criteria were not met
and/or see below

XII. QUANTITATION LIMITS

A. Dilution performed

(R)

041451515	T B. U. 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	
SAMPLE ID	DILUTION FACTOR	REASONS FOR DILUTION
Dilution was per	formed on samples by a	a factor of 2.29 except the following:
1502113AR1-	24.2	Analyte over the calibration range.
01A		
1502113AR1- 04A	12.4	2-Propanol over the calibration range after dilution. Result qualified as an estimated value (J) and qualified (E) by the laboratory.
V.		

B.	Percent Solids
	List samples which have ≤ 50 % solids
Actions	: If the % solids of a soil sample is 10-50%, estimate positive results (J) and nondetects (UJ)
	If the % solids of a soil sample is < 10%, estimate positive results (J) and reject nondetects